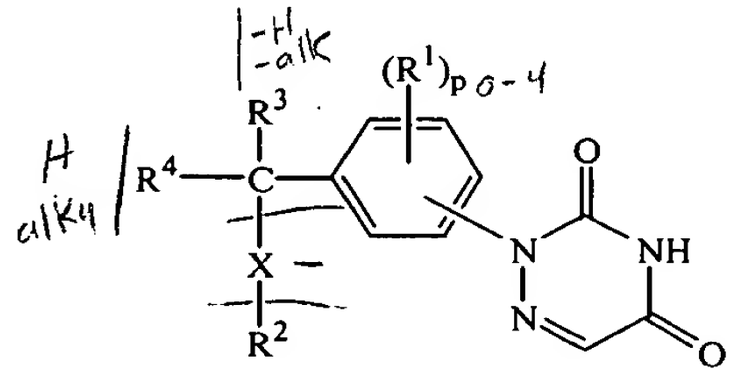
Note: Applicant uses:

- ~~*Cross-out text*~~ *to indicate deletions*
- *Underline text to indicate additions*

Claims:

1. (Currently Amended) A compound having the formula



(I)

a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein:

p represents an integer being 0, 1, 2, 3 or 4;

X represents O, S, NR^5 or a direct bond;

Y represents O, S, NR^5, or $S(O)_2$;

each R^1 independently represents C_{1-6}alkyl, halo, polyhaloC_{1-6}alkyl, hydroxy, mercapto, C_{1-6}alkyloxy, C_{1-6}alkylthio, C_{1-6}alkylcarbonyloxy, aryl, cyano, nitro, Het^3, R^6, NR^7R^8 or C_{1-4}alkyl substituted with Het^3, R^6 or NR^7R^8;

R^2 represents Het^1, C_{3-7}cycloalkyl, C_{1-6}alkyl or C_{1-6}alkyl substituted with one or two substituents selected from hydroxy, cyano, amino, mono- or di(C_{1-4}alkyl)amino, C_{1-6}alkyloxy, C_{1-6}alkylsulfonyloxy, C_{1-6}alkyloxycarbonyl, C_{3-7}cycloalkyl, aryl, aryloxy, arylthio, Het^1, Het^1oxy and Het^1thio; and if X is O, S or NR^5, then R^2 may also represent aminocarbonyl, aminothiocarbonyl, C_{1-4}alkylcarbonyl, C_{1-4}alkylthiocarbonyl, arylcarbonyl, arylthiocarbonyl, Het^1carbonyl or Het^1thiocarbonyl;

R^3 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl;

R^4 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl; or

R^3 and R^4 taken together form a C_{2-6}alkanediyl;

R^5 represents hydrogen or C_{1-4}alkyl;

each R^6 independently represents C_{1-6}alkylsulfonyl, aminosulfonyl, mono- or di(C_{1-4}alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl, polyhaloC_{1-6}alkylsulfonyl, C_{1-6}alkylsulfinyl, phenylC_{1-4}alkylsulfonyl, piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C_{1-4}alkyl-*N*-piperidinylaminosulfonyl or mono-or di(C_{1-4}alkyl)aminoC_{1-4}alkylsulfonyl;

each R^7 and each R^8 are independently selected from **the group consisting of:** hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

R^9 and R^{10} are each independently selected from **the group consisting of:** hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

each R^{11} independently being selected from **the group consisting of:** hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C_{1-4}alkyloxy, formyl, trihaloC_{1-4}alkylsulfonyloxy, R^6, NR^7R^8, C(=O)NR^7R^8, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, aryl, aryloxy, arylcarbonyl, C_{3-7}cycloalkyl, C_{3-7}cycloalkyloxy, phthalimide-2-yl, Het^3 and C(=O)Het^3;

R^{12} and R^{13} are each independently selected from **the group consisting of:** hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, phenylcarbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, C_{3-}

$_{7}$cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14} and R^{6};

each R^{14} independently represents hydrogen, C_{1-4}alkyl, C_{3-7}cycloalkyl, aminocarbonylmethylene or mono-or di(C_{1-4}alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each independently selected from nitro, azido, cyano, halo, hydroxy, C_{1-4}alkyl, C_{3-7}cycloalkyl, C_{1-4}alkyloxy, formyl, polyhaloC_{1-4}alkyl, $NR^{9}R^{10}$, $C(=O)NR^{9}R^{10}$, C(=O)-O-R^{14}, R^{6}, -O-R^{6}, phenyl, Het^{3}, C(=O)Het^{3} and C_{1-4}alkyl substituted with hydroxy, C_{1-4}alkyloxy, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^{3} or $NR^{9}R^{10}$;

Het^{1} represents a heterocycle selected from **the group consisting of:** pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiomorpholinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^{2}, R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from Het^{2} and R^{11};

provided Het^{1} is other than 2-substituted-pyridin-5-yl;

Het^{2} represents a heterocycle selected from **the group consisting of:** pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl,

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isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^4, R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from Het^4 and R^{11};

Het^3 represents a monocyclic heterocycle selected **the group consisting of:** from pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiomorpholinyl and tetrahydropyranyl; wherein said monocyclic heterocycles each independently may optionally be substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy, C_{1-4}alkyl, C_{1-4}alkyloxy, C_{1-4}alkylcarbonyl, piperidinyl, $NR^{12}R^{13}$, C(=O)-O-R^{14}, R^6 and C_{1-4}alkyl substituted with one or two substituents independently selected from hydroxy, C_{1-4}alkyloxy, phenyl, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, R^6 and $NR^{12}R^{13}$;

Het^4 represents a monocyclic heterocycle selected from **the group consisting of:** pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl.

2. (Currently Amended) A compound as claimed in claim 1 wherein:

each R^7 and each R^8 are independently selected from **the group consisting of:** hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 and R^6;

R^9 and R^{10} are each independently selected from **the group consisting of:** hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl,

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Het^{3}carbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^{3}aminocarbonyl, Het^{3}aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 and R^6;

R^{11} is being selected from **the group consisting of:** hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C_{1-4}alkyloxy, formyl, trihaloC_{1-4}alkylsulfonyloxy, R^6, NR^{7}R^8, C(=O)NR^{7}R^8, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, aryl, aryloxy, arylcarbonyl, C_{3-7}cycloalkyl, C_{3-7}cycloalkyloxy, phthalimide-2-yl, Het^3, Het^4 and C(=O)Het^3; and

Het^2 represents a heterocycle selected from **the group consisting of:** pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from R^{11}.

3-9. (Canceled)

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10. (Currently Amended) A compound **of the formula:** ~~**as claimed in claim 1 wherein the compound is**~~

2-[3,5-dichloro-4-[1-methyl-1-(4-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

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2-[3,5-dichloro-4-[1-[4-(3-chlorophenyl)-5-methyl-2-thiazolyl]-1-methylethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,2,4-oxadiazol-3-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-(4,5-diphenyl-2-thiazolyl)-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methylphenyl)-1,2,4-oxadiazol-3-yl]ethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-(4-methyl-5-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(3-pyridinyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(phenylmethyl)-2-thiazolyl]ethyl]-phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[5-(4-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[4-(3-thienyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-[4-(2-furanyl)-2-thiazolyl]-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[5-(3-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;

2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methyl-3-pyridinyl)-1,2, 4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; **or**

2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,3,4-oxadiazol-2-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; **or** a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof.

11. (Previously Amended) A composition comprising a pharmaceutically acceptable carrier and, as active ingredient, a therapeutically effective amount of a compound as claimed in claim 1.

12. through 18. (Previously Canceled)

19. through 55. (Currently Canceled)

55 D2 56. (New) A compound as claimed in claim 1 provided that in those compounds wherein X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

6 57. (New) A compound as claimed in claim 2 provided that in those compounds wherein X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

7 58. (New) A compound as claimed in claim 1 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

8 59. (New) A compound as claimed in claim 2 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

9 5 60. (New) A compound as claimed in claim 56 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

10 6 61. (New) A compound as claimed in claim 57 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

11 62. (New) A compound as claimed in claim 1 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~63.~~ A compound as claimed in claim 2 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~64.~~ A compound as claimed in claim ~~56~~ 5 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~65.~~ A compound as claimed in claim ~~57~~ 6 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~66.~~ A compound as claimed in claim ~~58~~ 7 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~67~~. A compound as claimed in claim ~~59~~ wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~68~~. A compound as claimed in claim ~~60~~ wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~69~~. A compound as claimed in claim ~~61~~ wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

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~~70~~. (New) A compound as claimed in claim 1 wherein R^3 and R^4 are both methyl and $-X-R^2$ is Het^1.

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~~71~~. (New) A compound as claimed in claim 2 wherein R^3 and R^4 are both methyl and $-X-R^2$ is Het^1.

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72. (New) A compound as claimed in claim 56 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

73. (New) A compound as claimed in claim 57 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

74. (New) A compound as claimed in claim 58 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

75. (New) A compound as claimed in claim 59 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

76. (New) A compound as claimed in claim 60 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

77. (New) A compound as claimed in claim 61 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

78. (New) A compound as claimed in claim 62 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

79. (New) A compound as claimed in claim 63 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

80. (New) A compound as claimed in claim 64 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

81. (New) A compound as claimed in claim 65 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

82. (New) A compound as claimed in claim 66 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

83. (New) A compound as claimed in claim 67 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

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~~84~~ 33. (New) A compound as claimed in claim ~~68~~ 17 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

~~85~~ 34. (New) A compound as claimed in claim ~~69~~ 18 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

~~86~~ 35. (New) A compound as claimed in claim ~~71~~ 1 wherein p is 1 or 2 and each R^1 is chloro.

~~87~~ 36. (New) A compound as claimed in claim ~~72~~ 2 wherein p is 1 or 2 and each R^1 is chloro.

~~88~~ 37. (New) A compound as claimed in claim ~~56~~ 5 wherein p is 1 or 2 and each R^1 is chloro.

~~89~~ 38. (New) A compound as claimed in claim ~~57~~ 6 wherein p is 1 or 2 and each R^1 is chloro.

~~90~~ 39. (New) A compound as claimed in claim ~~58~~ 7 wherein p is 1 or 2 and each R^1 is chloro.

~~91~~ 40. (New) A compound as claimed in claim ~~59~~ 8 wherein p is 1 or 2 and each R^1 is chloro.

~~92~~ 41. (New) A compound as claimed in claim ~~60~~ 9 wherein p is 1 or 2 and each R^1 is chloro.

~~93~~ 42. (New) A compound as claimed in claim ~~61~~ 10 wherein p is 1 or 2 and each R^1 is chloro.

~~94~~ 43. (New) A compound as claimed in claim ~~62~~ 11 wherein p is 1 or 2 and each R^1 is chloro.

95. (New) A compound as claimed in claim 12 wherein p is 1 or 2 and each R^1 is chloro.

96. (New) A compound as claimed in claim 13 wherein p is 1 or 2 and each R^1 is chloro.

97. (New) A compound as claimed in claim 14 wherein p is 1 or 2 and each R^1 is chloro.

98. (New) A compound as claimed in claim 15 wherein p is 1 or 2 and each R^1 is chloro.

99. (New) A compound as claimed in claim 16 wherein p is 1 or 2 and each R^1 is chloro.

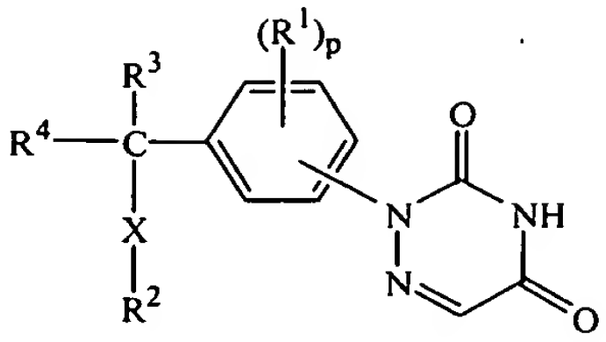
100. (New) A compound as claimed in claim 17 wherein p is 1 or 2 and each R^1 is chloro.

101. (New) A compound as claimed in claim 18 wherein p is 1 or 2 and each R^1 is chloro.

102. (New) A compound as claimed in claim 1 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

103. (New) A compound as claimed in claim 2 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

104. (New) A compound having the formula



(I)

a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein :

p represents an integer being 0, 1 or 2;

X represents O, S, NR^5 or a direct bond;

Y represents O, S, NR^5, or $S(O)_2$;

each R^1 independently represents chloro or trifluoromethyl;

R^2 represents Het^1 or C_{1-6}alkyl substituted with one or two substituents selected from hydroxy, cyano, amino, mono- or di(C_{1-4}alkyl)amino, C_{1-6}alkyloxy, C_{1-6}alkylsulfonyloxy, C_{1-6}alkyloxycarbonyl, C_{3-7}cycloalkyl, aryl, aryloxy, arylthio, Het^1, Het^1oxy and Het^1thio; and if X is O, S or NR^5, then R^2 may also represent aminocarbonyl, aminothiocarbonyl, C_{1-4}alkylcarbonyl, C_{1-4}alkylthiocarbonyl, arylcarbonyl, arylthiocarbonyl, Het^1carbonyl or Het^1thiocarbonyl;

R^3 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl;

R^4 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl; or R^3 and R^4 taken together form a C_{2-6}alkanediyl;

R^5 represents hydrogen or C_{1-4}alkyl;

each R^6 independently represents C_{1-6}alkylsulfonyl, aminosulfonyl, mono- or di(C_{1-4}alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl, polyhaloC_{1-6}alkylsulfonyl, C_{1-6}alkylsulfinyl, phenylC_{1-4}alkylsulfonyl, piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C_{1-4}alkyl-*N*-piperidinylaminosulfonyl or mono-or di(C_{1-4}alkyl)aminoC_{1-4}alkylsulfonyl;

each R^7 and each R^8 are independently selected from hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl,

C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

R^9 and R^{10} are each independently selected from hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

each R^{11} independently being selected from hydroxy, cyano, nitro, halo, C_{1-4}alkyloxy, formyl, NR^7R^8, C(=O)NR^7R^8, -C(=O)-O-R^{14}, aryl, arylcarbonyl, Het^3 and C(=O)Het^3;

each R^{14} independently represents hydrogen, C_{1-4}alkyl, C_{3-7}cycloalkyl, aminocarbonylmethylene or mono-or di(C_{1-4}alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each independently selected from nitro, azido, cyano, halo, hydroxy, C_{1-4}alkyl, C_{3-7}cycloalkyl, C_{1-4}alkyloxy, formyl, polyhaloC_{1-4}alkyl, NR^9R^{10}, C(=O)NR^9R^{10}, C(=O)-O-R^{14}, R^6, -O-R^6, phenyl, Het^3, C(=O)Het^3 and C_{1-4}alkyl substituted with hydroxy, C_{1-4}alkyloxy, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 or NR^9R^{10};

Het^1 represents a heterocycle selected from pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}; provided Het^1 is other than 2-substituted-pyridin-5-yl;

Het^2 represents a heterocycle selected from furanyl, thienyl, pyridinyl or benzothienyl, wherein said aromatic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^4, R^{11} and C_{1-4}alkyl optionally substituted with R^{11};

Het^3 represents a monocyclic heterocycle selected from piperidinyl, piperazinyl, morpholinyl and tetrahydropyranyl each independently and optionally substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy, C_{1-4}alkyl,

$C(=O)$-O-R^{14}, C_{1-4}alkylcarbonyl, R^6, piperidinyl and C_{1-4}alkyl substituted with one or two substituents independently selected from hydroxy, C_{1-4}alkyloxy, $C(=O)$-O-R^{14} and phenyl; Het^4 represents a monocyclic heterocycle selected from thienyl or pyridinyl.

D2 54 105. (New) A compound as claimed in claim 53 104, wherein when X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl, then R^2 is not substituted in the 6 position with an optionally substituted alkyl or acyl group.

55 106. (New) A compound as claimed in claim 53 104 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

56 107. (New) A compound as claimed in claim 53 104 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

57 108. (New) A compound as claimed in claim 53 104 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

58 109. (New) A compound as claimed in claim 53 104 wherein p is 1 or 2 and each R^1 is chloro.

59 110. (New) A compound as claimed in claim 53 104 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

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